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AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listing of claims in the application.

1. (Presently amended) A compound of structural formula 1:

$$(R^2)_n$$
 $(R^2)_n$
 $(R^2)_n$
 $(R^2)_n$
 $(R^2)_n$

I

a pharmaceutically acceptable salt or a stereoisomer thereof,

wherein:

a and b are each independently chosen from a double bond and a single bond;

X is hydrogen or halogen;

when a is a single bond, Y and Z are each independently chosen from hydrogen, C₁₋₄ alkyl, and halogen, or Y and Z, together with the carbon atom to which they are attached, form a cyclopropyl group;

when a is a double bond, Y is chosen from hydrogen, C_{1-4} alkyl, and halogen;

n is 0, 1, 2, or 3;

- U, V, W, and D are each independently chosen from CH, N, and S, and O, provided that at least one of U, V, W, and D is chosen from N and S, and O, and further provided that when one of U, V, W, and D is S or O, then the other ring members are independently chosen from N and CH;
- R¹ is chosen from hydrogen, CF₃, carbonyl(C₁₋₃ alkyl), hydroxyl, C₁₋₄ alkoxy, halogen, C₁₋₃ alkyl, hydroxymethyl, and (C₀₋₆ alkyl)₂amino, wherein said alkyl and alkoxy are each optionally substituted with one to seven fluorine atoms;

R² is chosen from:

halogen,
(carbonyl)₀₋₁C₁₋₁₀ alkyl,
(carbonyl)₀₋₁C₂₋₁₀ alkenyl,

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(carbonyl)₀₋₁C₂₋₁₀ alkynyl, C₁₋₁₀ alkenylamino,

(carbonyl)₀₋₁aryl C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl,

(C3-8)heterocyclyl C0-10 alkyl,

C3-8 heterocycloalkyl C0-10 alkyl,

C₁₋₄acylamino C₀₋₁₀ alkyl,

C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

di-(C₁₋₁₀ alkyl)amino C₀₋₁₀ alkyl,

arylC₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

(arylC0-10 alkyl)2amino C0-10 alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

(C3-8 cycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

(C3-8 heterocyclyl C0-10 alkyl)2amino C0-10 alkyl,

(C3-8 heterocycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂aminocarbonylamino,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonylamino,

C₀₋₁₀ alkyl aminocarbonylamino,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl aminocarbonylamino,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl C₀₋₁₀ alkyl,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C3-8 heterocycloalkyl C0-10 alkyl carbonylamino C0-10 alkyl,

aryl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

amino C₀₋₁₀ alkyl carbimidoylC₀₋₁₀ alkylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl,

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(aryl C<sub>1-10</sub> alkyl)<sub>1-2</sub>aminocarbonyl,
C<sub>1-10</sub> alkoxy (carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkyl,
C<sub>0-10</sub> alkylcarboxy C<sub>0-10</sub> alkylamino,
carboxy C<sub>0-10</sub> alkyl,
carboxy aryl,
carboxy C<sub>3-8</sub> cycloalkyl,
carboxy C3-8 heterocyclyl,
carboxy C3-8 heterocycloalkyl,
C<sub>1-10</sub> alkoxy,
C<sub>1-10</sub>alkyloxy C<sub>0-10</sub>alkyl,
aryloxy,
C<sub>3-8</sub> cycloalkyloxy,
C<sub>3-8</sub> heterocyclyloxy,
C<sub>3-8</sub> heterocycloalkyloxy,
C<sub>1-10</sub> alkylcarbonyloxy,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylcarbonyloxy,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylcarbonyloxy,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylcarbonyloxy,
aryl C<sub>0-10</sub> alkylcarbonyloxy,
C<sub>1-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
C3-8 heterocyclyl C0-10 alkyloxy(carbonyl)0-1C0-10 alkylamino,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
aryl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
(C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyloxy,
(aryl C<sub>0-10</sub> alkyl)<sub>1-2</sub>aminocarbonyloxy,
(C3-8 heterocyclyl C0-10 alkyl)1-2aminocarbonyloxy,
(C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyl)<sub>1-2</sub>aminocarbonyloxy,
(C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub>alkyl)<sub>1-2</sub>aminocarbonyloxy,
hydroxy C<sub>0-10</sub>alkyl,
hydroxycarbonylC0-10alkoxy,
hydroxycarbonylC0-10alkyloxy,
C<sub>1-10</sub> alkylthio,
C<sub>1-10</sub> alkylsulfinyl,
aryl C<sub>0-10</sub> alkylsulfinyl,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylsulfinyl,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylsulfinyl,
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C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylsulfinyl,
          C<sub>1-10</sub> alkylsulfonyl,
          aryl C<sub>0-10</sub> alkylsulfonyl,
          C3-8 heterocyclyl C0-10 alkylsulfonyl,
          C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylsulfonyl,
          C3-8 cycloalkyl C0-10 alkylsulfonyl,
          C<sub>1-10</sub> alkylsulfonylamino,
          aryl C<sub>1-10</sub> alkylsulfonylamino,
          C3-8 heterocyclyl C1-10 alkylsulfonylamino,
          C<sub>3-8</sub> heterocycloalkyl C<sub>1-10</sub> alkylsulfonylamino,
          C3-8 cycloalkyl C1-10 alkylsulfonylamino,
          cyano,
           nitro,
           perfluoroC<sub>1</sub>-6alkyl, and
           perfluoroC<sub>1-6</sub>alkoxy, and
wherein R<sup>2</sup> is optionally substituted with at least one substituent, R<sup>3</sup>, chosen from:
           halogen,
           (carbonyl)0-1C1-10 alkyl,
           (carbonyl)0-1C2-10 alkenyl,
           (carbonyl)0-1C2-10 alkynyl,
           (carbonyl)<sub>0-1</sub>aryl C<sub>0-10</sub> alkyl,
           C3-8 cycloalkyl C0-10 alkyl,
           (C3-8)heterocyclyl C0-10 alkyl,
           (C3-8)heterocycloalkyl C0-10 alkyl,
           C<sub>1</sub>-4acylamino C<sub>0-10</sub> alkyl,
           C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           di-(C1-10 alkyl)amino C0-10 alkyl,
           arylC0-10 alkylamino C0-10 alkyl,
           (arylC<sub>0-10</sub> alkyl)<sub>2</sub>amino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>0-10</sub> alkyl carbimidoylC<sub>0-10</sub> alkyl,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyl,
           C<sub>1-10</sub> alkoxy (carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkyl,
           C<sub>1-10</sub>alkyloxy C<sub>0-10</sub>alkyl,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyloxy,
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hydroxycarbonylC0-10alkoxy,

(C₁₋₁₀ alkyl)2aminocarbonyloxy,

(aryl C₀₋₁₀ alkyl)₁₋₂aminocarbonyloxy,

hydroxy C₀₋₁₀alkyl,

C₁₋₁₀ alkylsulfonyl,

C₁₋₁₀ alkylsulfonylamino,

aryl C₁₋₁₀ alkylsulfonylamino,

C₃₋₈ heterocyclyl C₁₋₁₀ alkylsulfonylamino,

C₃₋₈ heterocycloalkyl C₁₋₁₀ alkylsulfonylamino,

C3-8 cycloalkyl C1-10 alkylsulfonylamino,

cyano,

nitro,

perfluoroC₁₋₆alkyl, and

perfluoroC₁₋₆alkoxy,

wherein R³ is optionally substituted with one or more groups chosen from hydrogen, OH, (C₁-6)alkoxy, halogen, CO₂H, CN, O(C=O)C₁-C₆ alkyl, NO₂, trifluoromethoxy, trifluoroethoxy, -O₍₀₋₁₎(C₁₋₁₀)perfluoroalkyl, and NH₂.

- 2. (Original) A compound according to Claim 1, wherein X is fluorine.
- 3. (Original) A compound according to Claim 1, wherein X is hydrogen.
- 4. (Original) A compound according to Claim 1, wherein a is a single bond and b is a double bond.
- (Original) A compound according to Claim 1 and of structural formula II, wherein:

$$(R^2)_n$$
 $(R^2)_n$
 $(R^2)_n$

a pharmaceutically acceptable salt or a stereoisomer thereof,

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wherein:

a and b are each independently chosen from a double bond and a single bond;

n is 0, 1, 2, or 3;

X is hydrogen or halogen;

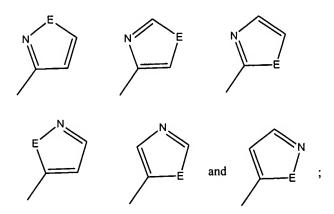
when a is a single bond, Y and Z are each independently chosen from hydrogen, C₁₋₄ alkyl, and

halogen, or Y and Z, together with the carbon atom to which they are attached, form a cyclopropyl group;

when a is a double bond, Y is chosen from hydrogen, C₁₋₄ alkyl, and halogen;



is chosen from:



E is S or O;

R¹ is chosen from: hydrogen, CF₃, carbonyl(C₁₋₃ alkyl), hydroxyl, C₁₋₄ alkoxy, halogen, C₁₋₃ alkyl, hydroxymethyl, and (C₀₋₆ alkyl)₂amino, wherein said alkyl and alkoxy are each optionally substituted with one to seven fluorine atoms;

R² is chosen from:

halogen,

(carbonyl)0-1C1-10 alkyl,

(carbonyl)₀₋₁C₂₋₁₀ alkenyl,

(carbonyl)0-1C2-10 alkynyl,

C₁₋₁₀ alkenylamino,

(carbonyl)0-1aryl C0-10 alkyl,

C3-8 cycloalkyl C0-10 alkyl,

(C3-8)heterocyclyl C0-10 alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl,

C₁₋₄acylamino C₀₋₁₀ alkyl,

C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

di-(C1-10 alkyl)amino C0-10 alkyl,

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arylC0-10 alkylamino C0-10 alkyl,

(arylC₀₋₁₀ alkyl)2amino C₀₋₁₀ alkyl,

C₁₋₁₀ alkoxy (carbonyl)₀₋₁C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

(C3-8 cycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

(C3-8 heterocyclyl C0-10 alkyl)2amino C0-10 alkyl,

(C3-8 heterocycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂ aminocarbonylamino,

(aryl C1-10 alkyl)1-2aminocarbonylamino,

C₀₋₁₀ alkyl aminocarbonylamino,

C3-8 heterocyclyl C0-10 alkyl aminocarbonylamino,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl C₀₋₁₀ alkyl,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C3-8 cycloalkyl C0-10 alkyl aminocarbonyl C0-10 alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

amino C₀₋₁₀ alkyl carbimidoylC₀₋₁₀ alkylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl,

(aryl C₁₋₁₀ alkyl)₁₋₂ aminocarbonyl,

C₁₋₁₀ alkoxy (carbonyl)₀₋₁C₀₋₁₀ alkyl,

C₀₋₁₀ alkylcarboxy C₀₋₁₀ alkylamino,

carboxy C₀₋₁₀ alkyl,

carboxy aryl,

carboxy C₃₋₈ cycloalkyl,

carboxy C3.8 heterocyclyl,

carboxy C3-8 heterocycloalkyl,

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C₁₋₁₀ alkoxy,

C₁₋₁₀alkyloxy C₀₋₁₀alkyl,

aryloxy,

C3-8 cycloalkyloxy,

C3-8 heterocyclyloxy,

C3-8 heterocycloalkyloxy,

C₁₋₁₀ alkylcarbonyloxy,

C3-8 heterocyclyl C0-10 alkylcarbonyloxy,

C3-8 heterocycloalkyl C0-10 alkylcarbonyloxy,

C₃₋₈ cycloalkyl C₀₋₁₀ alkylcarbonyloxy,

aryl C₀₋₁₀ alkylcarbonyloxy,

C₁₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C3-8 cycloalkyl C0-10 alkyloxy(carbonyl)0-1C0-10 alkylamino,

aryl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

(C₁₋₁₀ alkyl)2aminocarbonyloxy,

(aryl C₀₋₁₀ alkyl)₁₋₂aminocarbonyloxy,

(C3-8 heterocyclyl C0-10 alkyl)1-2aminocarbonyloxy,

(C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl)₁₋₂aminocarbonyloxy,

(C3-8 cycloalkyl C0-10alkyl)1-2aminocarbonyloxy,

hydroxy C₀₋₁₀alkyl,

hydroxycarbonylC0-10alkoxy,

hydroxycarbonylC₀₋₁0alkyloxy,

C₁₋₁₀ alkylthio,

C₁₋₁₀ alkylsulfinyl,

aryl C₀₋₁₀ alkylsulfinyl,

C3-8 heterocyclyl C0-10 alkylsulfinyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkylsulfinyl,

C3-8 cycloalkyl C0-10 alkylsulfinyl,

C₁₋₁₀ alkylsulfonyl,

aryl C₀₋₁₀ alkylsulfonyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkylsulfonyl,

C3-8 heterocycloalkyl C0-10 alkylsulfonyl,

C3-8 cycloalkyl C0-10 alkylsulfonyl,

C₁₋₁₀ alkylsulfonylamino,

aryl C1-10 alkylsulfonylamino,

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C3-8 heterocyclyl C1-10 alkylsulfonylamino,
           C<sub>3-8</sub> heterocycloalkyl C<sub>1-10</sub> alkylsulfonylamino,
           C<sub>3-8</sub> cycloalkyl C<sub>1-10</sub> alkylsulfonylamino,
           cyano,
           nitro,
           perfluoroC<sub>1-6</sub>alkyl, and
           perfluoroC<sub>1-6</sub>alkoxy, and
wherein R<sup>2</sup> is optionally substituted with at least one substituent R<sup>3</sup>;
R<sup>3</sup> is chosen from:
          halogen,
           (carbonyl)<sub>0-1</sub>C<sub>1-10</sub> alkyl,
           (carbonyl)0-1C2-10 alkenyl,
           (carbonyl)0-1C2-10 alkynyl,
           (carbonyl)0-1aryl C0-10 alkyl,
           C3-8 cycloalkyl C0-10 alkyl,
           (C3-8)heterocyclyl C0-10 alkyl,
           (C3-8)heterocycloalkyl C0-10 alkyl,
           C<sub>1</sub>-4acylamino C<sub>0</sub>-10 alkyl,
           C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           di-(C<sub>1-10</sub> alkyl)amino C<sub>0-10</sub> alkyl,
           arylC<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           (arylC0-10 alkyl)2amino C0-10 alkyl,
           C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>0-10</sub> alkyl carbimidoylC<sub>0-10</sub> alkyl,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyl,
           C1-10 alkoxy (carbonyl)0-1C0-10 alkyl,
           C<sub>1-10</sub>alkyloxy C<sub>0-10</sub>alkyl,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyloxy,
           hydroxycarbonylC0-10alkoxy,
           (C1-10 alkyl)2aminocarbonyloxy,
           (aryl C<sub>0-10</sub> alkyl)<sub>1-2</sub>aminocarbonyloxy,
           hydroxy C<sub>0-10</sub>alkyl,
           C<sub>1-10</sub> alkylsulfonyl,
           C<sub>1-10</sub> alkylsulfonylamino,
           aryl C<sub>1-10</sub> alkylsulfonylamino,
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C3-8 heterocyclyl C1-10 alkylsulfonylamino,
C3-8 heterocycloalkyl C1-10 alkylsulfonylamino,
C3-8 cycloalkyl C1-10 alkylsulfonylamino,
cyano,
nitro,
perfluoroC1-6alkyl, and
perfluoroC1-6alkoxy,
wherein R3 is optionally substituted with one or more groups chosen from hydrogen, OH, (C1-6)alkoxy, halogen, CO2H, CN, O(C=O)C1-C6 alkyl, NO2, trifluoromethoxy, trifluoroethoxy,
-O(0-1)(C1-10)perfluoroalkyl, and NH2.
```

- 6. (Original) A compound according to Claim 5, wherein R¹ is chosen from: hydrogen, CF₃, hydroxyl, and C₁₋₃ alkyl optionally substituted with one to seven fluorine atoms.
- 7. (Original) A compound according to Claim 6, wherein R¹ is chosen from: hydrogen and C₁₋₃ alkyl.
 - 8. (Original) A compound according to Claim 7, wherein R¹ is methyl.
 - 9. (Original) A compound according to Claim 8, wherein R² is chosen from:

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halogen,
(carbonyl)<sub>0-1</sub>C<sub>1-10</sub> alkyl,
(carbonyl)0-1C2-10 alkenyl,
(carbonyl)0-1C2-10 alkynyl,
C<sub>1-10</sub> alkenylamino,
(carbonyl)0-1 aryl C0-10 alkyl,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkyl,
(C<sub>3-8</sub>)heterocyclyl C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyl,
C<sub>1-4</sub> acylamino C<sub>0-10</sub> alkyl,
C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
di-(C<sub>1-10</sub> alkyl)amino C<sub>0-10</sub> alkyl,
arylC<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
(arylC<sub>0-10</sub> alkyl)<sub>2</sub>amino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
(C3-8 cycloalkyl C0-10 alkyl)2amino C0-10 alkyl,
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(C3-8 heterocyclyl C0-10 alkyl)2amino C0-10 alkyl,
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(C3-8 heterocycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂aminocarbonylamino,

(aryl C1-10 alkyl)1-2aminocarbonylamino,

C₀₋₁₀ alkyl aminocarbonylamino,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl aminocarbonylamino,

C3-8 heterocycloalkyl C0-10 alkyl aminocarbonylamino,

C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

amino C₀₋₁₀ alkyl carbimidoylC₀₋₁₀ alkylamino,

C₀₋₁₀ alkylcarboxy C₀₋₁₀ alkylamino,

C₁₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

aryl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C₁₋₁₀ alkylsulfonylamino,

aryl C₁₋₁₀ alkylsulfonylamino,

C₃₋₈ heterocyclyl C₁₋₁₀ alkylsulfonylamino,

C3-8 heterocycloalkyl C1-10 alkylsulfonylamino,

C₃₋₈ cycloalkyl C₁₋₁₀ alkylsulfonylamino,

cyano,

nitro,

perfluoroC₁₋₆alkyl, and

perfluoroC1-6alkoxy, and

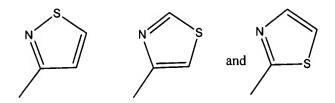
wherein R² is optionally substituted with at least one substituent R³.

10. (Original) A compound according to Claim 9, wherein E is S.

11. (Original) A compound according to Claim 10, wherein



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- 12. (Original) A compound according to Claim 11, wherein b is a double bond.
- 13. (Original) A compound according to Claim 12, wherein a is a single bond and b is a double bond.
 - 14. (Original) A compound according to Claim 1 and of structural formula III,

$$(R^2)_n$$
 $(R^2)_n$
 $(R^2$

a pharmaceutically acceptable salt or a stereoisomer thereof, wherein:

X is hydrogen or halogen;

n is 0, 1, 2, or 3;

Y and Z are each independently chosen from hydrogen, C₁₋₄ alkyl, and halogen, or Y and Z, together with the carbon atom to which they are attached, form a cyclopropyl group;

U, V, W, and D are each independently chosen from N and CH, provided that at least one of U, V, W, and D is CH;

R² is chosen from:

halogen,

(carbonyl)0-1C1-10 alkyl,

(carbonyl)0-1C2-10 alkenyl,

(carbonyl)₀₋₁C₂₋₁₀ alkynyl,

C₁₋₁₀ alkenylamino,

(carbonyl)0-1aryl C0-10 alkyl,

C3-8 cycloalkyl C0-10 alkyl,

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(C3-8)heterocyclyl C0-10 alkyl,

C3-8 heterocycloalkyl C0-10 alkyl,

C₁-4acylamino C₀-10 alkyl,

C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

di-(C1-10 alkyl)amino C0-10 alkyl,

arylC0-10 alkylamino C0-10 alkyl,

(arylC0-10 alkyl)2amino C0-10 alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkylamino C₀₋₁₀ alkyl,

(C₃₋₈ cycloalkyl C₀₋₁₀ alkyl)₂amino C₀₋₁₀ alkyl,

(C₃₋₈ heterocyclyl C₀₋₁₀ alkyl)₂amino C₀₋₁₀ alkyl,

(C3-8 heterocycloalkyl C0-10 alkyl)2amino C0-10 alkyl,

C3-8 cycloalkyl C0-10 alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)2aminocarbonylamino,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonylamino,

C₀₋₁₀ alkyl aminocarbonylamino,

C3-8 heterocyclyl C0-10 alkyl aminocarbonylamino,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl C₀₋₁₀ alkyl,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl aminocarbonyl C₀₋₁₀ alkyl,

C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocyclyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

aryl C₀₋₁₀ alkyl carbonylamino C₀₋₁₀ alkyl,

amino C₀₋₁₀ alkyl carbimidoylC₀₋₁₀ alkylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyl,

(aryl C₁₋₁₀ alkyl)₁₋₂aminocarbonyl,

C₁₋₁₀ alkoxy (carbonyl)₀₋₁C₀₋₁₀ alkyl,

C₀₋₁₀ alkylcarboxy C₀₋₁₀ alkylamino,

carboxy C₀₋₁₀ alkyl,

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carboxy aryl,
carboxy C<sub>3-8</sub> cycloalkyl,
carboxy C<sub>3-8</sub> heterocyclyl,
carboxy C<sub>3-8</sub> heterocycloalkyl,
C<sub>1-10</sub> alkoxy,
```

C1-10alkyloxy C0-10alkyl,

aryloxy,

C₃₋₈ cycloalkyloxy,

C3-8 heterocyclyloxy,

C3-8 heterocycloalkyloxy,

C₁₋₁₀ alkylcarbonyloxy,

C₃₋₈ heterocyclyl C₀₋₁₀ alkylcarbonyloxy,

C3-8 heterocycloalkyl C0-10 alkylcarbonyloxy,

C3-8 cycloalkyl C0-10 alkylcarbonyloxy,

aryl C₀₋₁₀ alkylcarbonyloxy,

C₁₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

C3-8 heterocyclyl C0-10 alkyloxy(carbonyl)0-1C0-10 alkylamino,

C3-8 heterocycloalkyl C0-10 alkyloxy(carbonyl)0-1C0-10 alkylamino,

C3-8 cycloalkyl C0-10 alkyloxy(carbonyl)0-1C0-10 alkylamino,

aryl C₀₋₁₀ alkyloxy(carbonyl)₀₋₁C₀₋₁₀ alkylamino,

(C₁₋₁₀ alkyl)₂aminocarbonyloxy,

(aryl C₀₋₁₀ alkyl)₁₋₂aminocarbonyloxy,

(C3-8 heterocyclyl C0-10 alkyl)1-2aminocarbonyloxy,

(C₃₋₈ heterocycloalkyl C₀₋₁₀ alkyl)₁₋₂aminocarbonyloxy,

(C3-8 cycloalkyl C0-10alkyl)1-2aminocarbonyloxy,

hydroxy C₀₋₁₀alkyl,

hydroxycarbonylC0-10alkoxy,

hydroxycarbonylC0-10alkyloxy,

C₁₋₁₀ alkylthio,

C₁₋₁₀ alkylsulfinyl,

aryl C₀₋₁₀ alkylsulfinyl,

C3-8 heterocyclyl C0-10 alkylsulfinyl,

C3-8 heterocycloalkyl C0-10 alkylsulfinyl,

C₃₋₈ cycloalkyl C₀₋₁₀ alkylsulfinyl,

C₁₋₁₀ alkylsulfonyl,

aryl C₀₋₁₀ alkylsulfonyl,

C3-8 heterocyclyl C0-10 alkylsulfonyl,

```
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylsulfonyl,
           C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylsulfonyl,
           C<sub>1-10</sub> alkylsulfonylamino,
           aryl C<sub>1-10</sub> alkylsulfonylamino,
           C3-8 heterocyclyl C1-10 alkylsulfonylamino,
           C3-8 heterocycloalkyl C1-10 alkylsulfonylamino,
           C<sub>3-8</sub> cycloalkyl C<sub>1-10</sub> alkylsulfonylamino,
           cyano,
           nitro,
           perfluoroC1-6alkyl, and
           perfluoroC<sub>1</sub>-6alkoxy, and
wherein R<sup>2</sup> is optionally substituted with at least one substituent, R<sup>3</sup>, chosen from:
           halogen,
           (carbonyl)0-1C1-10 alkyl,
           (carbonyl)0-1C2-10 alkenyl,
           (carbonyl)<sub>0-1</sub>C<sub>2-10</sub> alkynyl,
           (carbonyl)0-1aryl C0-10 alkyl,
           C3-8 cycloalkyl C0-10 alkyl,
           (C<sub>3-8</sub>)heterocyclyl C<sub>0-10</sub> alkyl,
           (C<sub>3-8</sub>)heterocycloalkyl C<sub>0-10</sub> alkyl,
           C<sub>1-4</sub>acylamino C<sub>0-10</sub> alkyl,
           C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           di-(C<sub>1-10</sub> alkyl)amino C<sub>0-10</sub> alkyl,
           arylC<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           (arylC0-10 alkyl)2amino C0-10 alkyl,
           C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
           C<sub>0-10</sub> alkyl carbimidoylC<sub>0-10</sub> alkyl,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyl,
           C<sub>1-10</sub> alkoxy (carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkyl,
           C<sub>1-10</sub>alkyloxy C<sub>0-10</sub>alkyl,
           (C<sub>1-10</sub> alkyl)2aminocarbonyloxy,
           hydroxycarbonylC0-10alkoxy,
           (C<sub>1-10</sub> alkyl)<sub>2</sub>aminocarbonyloxy,
           (aryl C<sub>0-10</sub> alkyl)<sub>1-2</sub>aminocarbonyloxy,
           hydroxy C<sub>0-10</sub>alkyl,
```

```
C1-10 alkylsulfonylamino,
aryl C1-10 alkylsulfonylamino,
C3-8 heterocyclyl C1-10 alkylsulfonylamino,
C3-8 heterocycloalkyl C1-10 alkylsulfonylamino,
C3-8 cycloalkyl C1-10 alkylsulfonylamino,
C3-8 cycloalkyl C1-10 alkylsulfonylamino,
cyano,
nitro,
perfluoroC1-6alkyl, and
perfluoroC1-6alkoxy, and
wherein R³ is optionally substituted with one or more groups chosen from hydrogen, OH, (C1-6)alkoxy, halogen, CO2H, CN, O(C=O)C1-C6 alkyl, NO2, trifluoromethoxy, trifluoroethoxy,
-O(0-1)(C1-10)perfluoroalkyl, and NH2.
```

- 15. (Original) A compound according to Claim 14, wherein X is hydrogen.
- 16. (Original) A compound according to Claim 15, wherein R² is chosen from:

```
halogen,
(carbonyl)<sub>0-1</sub>C<sub>1-10</sub> alkyl,
(carbonyl)0-1C2-10 alkenyl,
(carbonyl)<sub>0-1</sub>aryl C<sub>0-10</sub> alkyl,
C3-8 cycloalkyl C0-10 alkyl,
(C3-8)heterocyclyl C0-10 alkyl,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyl,
C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
arylC<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkyl aminocarbonylamino,
(aryl C<sub>1-10</sub> alkyl)<sub>1-2</sub>aminocarbonylamino,
C<sub>0-10</sub> alkyl aminocarbonylamino,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkyl aminocarbonylamino,
C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyl aminocarbonylamino,
C<sub>0-10</sub> alkyl carbonylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkyl carbonylamino C<sub>0-10</sub> alkyl,
C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkyl carbonylamino C<sub>0-10</sub> alkyl,
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C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyl carbonylamino C<sub>0-10</sub> alkyl,
          aryl C<sub>0-10</sub> alkyl carbonylamino C<sub>0-10</sub> alkyl,
          C<sub>0-10</sub> alkylcarboxy C<sub>0-10</sub> alkylamino,
          C<sub>1-10</sub> alkoxy,
          C1-10alkyloxy C0-10alkyl,
          aryloxy,
          C3-8 cycloalkyloxy,
          C3-8 heterocyclyloxy,
          C3-8 heterocycloalkyloxy,
          C<sub>1-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
          C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
           C<sub>3-8</sub> heterocycloalkyl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
           C<sub>3-8</sub> cycloalkyl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
           aryl C<sub>0-10</sub> alkyloxy(carbonyl)<sub>0-1</sub>C<sub>0-10</sub> alkylamino,
           hydroxy C<sub>0-10</sub>alkyl,
           C<sub>1-10</sub> alkylthio,
           C<sub>1-10</sub> alkylsulfonyl,
           aryl C<sub>0-10</sub> alkylsulfonyl,
           C<sub>3-8</sub> heterocyclyl C<sub>0-10</sub> alkylsulfonyl,
           C3-8 heterocycloalkyl C0-10 alkylsulfonyl,
           C3-8 cycloalkyl C0-10 alkylsulfonyl,
           C<sub>1-10</sub> alkylsulfonylamino,
           aryl C<sub>1-10</sub> alkylsulfonylamino,
           C3-8 heterocyclyl C1-10 alkylsulfonylamino,
           C<sub>3-8</sub> heterocycloalkyl C<sub>1-10</sub> alkylsulfonylamino,
           C3-8 cycloalkyl C1-10 alkylsulfonylamino,
           cyano,
           nitro.
           perfluoroC1-6alkyl, and
           perfluoroC<sub>1</sub>-6alkoxy, and
wherein R<sup>2</sup> is optionally substituted with at least one substituent R<sup>3</sup>.
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- 17. (Original) A compound according to Claim 16, wherein at least two of U, V, W, and D are each N and provided that at least one of U, V, W, and D is CH
 - 18. (Original) A compound according to Claim 1, selected from:

 17β -[2-(butylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;

```
17\beta-[2-(anilino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyridin-2-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[(2-methylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-{[2-methyl(phenyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(4-fluorophenyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(benzylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(isopropylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyridin-3-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(2-fluorophenyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(methoxyethyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(2-piperid-1-ylethyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(t-butyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-{[2-(4-cyanophenyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(cyclohexyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-{2-[(pyridin-4-ylmethyl]amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(pyrimidin-2-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyridin-4-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-{2-[(cyclopropylmethyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(propylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(allylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(heptylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(\text{octylamino})-1,3-\text{thiazol-}4-\text{yl}]-4-\text{methyl-}4-\text{aza-}5\alpha-\text{androst-}1-\text{en-}3-\text{one};
17\beta-[2-(hexylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(5-\text{methyl-1},2,3-\text{thiadiazol-2-yl})\text{amino}]-1,3-\text{thiazol-4-yl}]-4-\text{methyl-4-aza-}5\alpha-\text{androst-1-en-3-}
        one;
17\beta-{[2-(methoxypropyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-{2-[(2-morphilin-1-ylethyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-\{2-[(2,2,2-\text{trifluoroethyl})\text{amino}]-1,3-\text{thiazol-4-yl}\}-4-\text{methyl-4-aza-}5\alpha-\text{androst-1-en-3-one};
17\beta-\{2-\lceil (pyridin-2-ylethyl]amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(guanidino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(1-methyl-1H-imidazole-5-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-
        one;
17\beta-[2-(acetamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(phenyl carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(thiephene-3-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(furan-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
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17\beta-[2-(pyrizine-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(pyridine-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(thiephene-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(pyridine-3-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(pyridine-4-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(1-t-butyl-3-methyl-1H-pyrazole-5-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-
        androst-1-en-3-one;
17β-[2-(1-methyl-proline-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(1-methyl-1H-imidazole-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-
17β-[2-(1H-imidazole-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(methanesulfonamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\(\alpha\)-androst-1-en-3-one;
17\beta-[2-(ethyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(isopropyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(2-fluoroethyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(t-butylcarbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one
17\beta-[2-(ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-pyridin-2ylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-cyclopropylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-cyclohexylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-cyclohexylmethylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(morpholine-4-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(piperizine-1-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[N'-isopropylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(pyridyl-3-ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[N'-(methylamino)ethethylureyl]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androstan-3-one;
17\beta-[2-(pyridin-2-yl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androstan-3-one;
17\beta-[2-(methyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyrid-3-yl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(ethyl acetyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(acetonitrily1)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(2-chlorophenyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(methyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(phenyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(3,5-dimethylpyrazol-1-yl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(aminoacetyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
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17\beta-[5-(ureyl)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[5-(N-methyl-ureyl)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[5-(N,N-dimethyl-ureyl)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5α-androst-1-en-3-one; and
pharmaceutically acceptable salts and stereoisomers thereof.
                 19.
                          (Original) A compound according to Claim 18, selected from:
17\beta-[2-(butylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(anilino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(pyridin-2-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[(2-methylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-methyl(phenyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-{[2-(4-fluorophenyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(benzylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(isopropylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyridin-3-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-{[2-(2-fluorophenyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(methoxyethyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(2-piperid-1-ylethyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(t-butyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{[2-(4-cyanophenyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(cyclohexyl)amino]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(pyridin-4-ylmethyl]amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyrimidin-2-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(pyridin-4-ylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-{2-[(cyclopropylmethyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(propylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(allylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(heptylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(\text{octylamino})-1,3-\text{thiazol-4-yl}]-4-\text{methyl-4-aza-}5\alpha-\text{androst-1-en-3-one};
17\beta-[2-(hexylamino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-\{2-[(5-methyl-1,2,3-thiadiazol-2-yl)amino]-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-
        one;
17β-{[2-(methoxypropyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-{2-[(2-morphilin-1-ylethyl)amino]-1,3-thiazol-4-yl}-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-\{2-[(2,2,2-\text{trifluoroethyl})\text{amino}]-1,3-\text{thiazol-4-yl}\}-4-\text{methyl-4-aza-}5\alpha-\text{androst-1-en-3-one};
17\beta-{2-[(pyridin-2-ylethyl]amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
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 17β -[5-(amino)-1,2,4-triazol-3-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;

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17\beta-[2-(amino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(guanidino)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(1-methyl-1H-imidazole-5-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-
        one;
17\beta-[2-(acetamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(phenyl carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(thiephene-3-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(furan-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(pyrizine-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(pyridine-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(thiephene-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(pyridine-3-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(pyridine-4-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(1-t-butyl-3-methyl-1H-pyrazole-5-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-
        androst-1-en-3-one;
17β-[2-(1-methyl-proline-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(1-methyl-1H-imidazole-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-
        one;
17β-[2-(1H-imidazole-2-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(methanesulfonamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(ethyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(isopropyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(2-fluoroethyl carbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(t-butylcarbamate)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one
17\beta-[2-(ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-pyridin-2ylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-cyclopropylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(N'-cyclohexylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17β-[2-(N'-cyclohexylmethylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(morpholine-4-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[2-(piperizine-1-carboxamido)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17β-[N'-isopropylureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-[2-(pyridyl-3-ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one;
17\beta-\{2-[N'-(methylamino)ethethylureyl]-1,3-thiazol-4-yl\}-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
17\beta-[2-(ureyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androstan-3-one;
17\beta-[2-(pyridin-2-yl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androstan-3-one;
17\beta-[2-(methyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5\alpha-androst-1-en-3-one;
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17β-[2-(pyrid-3-yl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one; 17β-[2-(ethyl acetyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one; 17β-[2-(acetonitrilyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one; 17β-[2-(2-chlorophenyl)-1,3-thiazol-4-yl]-4-methyl-4-aza-5α-androst-1-en-3-one; and pharmaceutically acceptable salts and stereoisomers thereof.
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20. (Original) A compound according to Claim 19, selected from: 17β -[2-(methyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[2-(phenyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[2-(3,5-dimethylpyrazol-1-yl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[2-(aminoacetyl)-1,3-imidazol-4-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[5-(amino)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[5-(ureyl)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; 17β -[5-(N,N-dimethyl-ureyl)- 1,2,4-triazol-3-yl]-4-methyl-4-aza-5 α -androst-1-en-3-one; and pharmaceutically acceptable salts and stereoisomers thereof.

21 to 23. (Cancelled)

24. (Presently amended) A method of treating a condition in a mammal which is eaused by androgen deficiency, which can be ameliorated by androgen replacement, or which can be increased by androgen replacement, which selected from weakened muscle tone, osteoporosis, osteopenia, glucocorticoid-induced osteoporosis, periodontal disease, bone fracture, bone damage following bone reconstructive surgery, sarcopenia, frailty, aging skin, male hypogonadism, postmenopausal symptoms in women, atherosclerosis, hypercholesterolemia, hyperlipidemia, obesity, aplastic anemia and other hematopoietic disorders, inflammatory arthritis and joint repair, HIV-wasting, prostate cancer, benign prostatic hyperplasia (BPH), abdominal adiposity, metabolic syndrome, type II diabetes, cancer cachexia, Alzheimer's disease, muscular dystrophies, cognitive decline, sexual dysfunction, sleep apnea, depression, premature ovarian failure, and autoimmune disease, comprising administering to the mammal in need of such treatment, a therapeutically effective amount of a compound according to Claim 1 or a pharmaceutically acceptable salt or a stereoisomer thereof.

25 to 30. (Cancelled)

31. (Original) A pharmaceutical composition comprising a therapeutically effective amount of a compound of Claim 1 and a pharmaceutically acceptable carrier.

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32. (Original) A composition of Claim 31, further comprising an active ingredient selected from:

- an estrogen or an estrogen derivative, alone or in combination with a progestin or progestin derivative,
- 2) a bisphosphonate,
- 3) an antiestrogen or a selective estrogen receptor modulator,
- 4) an ανβ3 integrin receptor antagonist,
- 5) a cathepsin K inhibitor,
- 6) an HMG-CoA reductase inhibitor,
- 7) an osteoclast vacuolar ATPase inhibitor,
- 8) an antagonist of VEGF binding to osteoclast receptors,
- 9) an activator of peroxisome proliferator-activated receptor γ ,
- 10) calcitonin,
- 11) a calcium receptor antagonist,
- 12) parathyroid hormone or analog thereof,
- 13) a growth hormone secretagogue,
- 14) human growth hormone,
- 15) insulin-like growth factor,
- 16) a p38 protein kinase inhibitor,
- 17) bone morphogenetic protein,
- 18) an inhibitor of BMP antagonism,
- 19) a prostaglandin derivative,
- 20) vitamin D or vitamin D derivative,
- 21) vitamin K or vitamin K derivative,
- 22) ipriflavone,
- 23) fluoride salts,
- 24) dietary calcium supplement, and
- 25) osteoprotegerin.
- 33. (Original) A composition of Claim 32, wherein said bisphosphonate is alendronate.

34 to 37. (Cancelled)

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38. A pharmaceutical composition made by combining a compound according to Claim 1 and a pharmaceutically acceptable carrier.

39. A process for making a pharmaceutical composition comprising combining a compound according to Claim 1 and a pharmaceutically acceptable carrier.

40 to 41. (Cancelled)